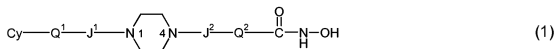


AMENDMENTS TO THE CLAIMS:

Please amend the claims as follows:

Claims 1-79. (Canceled).

80. (Currently Amended) A compound of the formula:



wherein:

the piperazin-1,4-diyl group is optionally substituted;

J¹ is independently a covalent bond or -C(=O)- ;

J² is independently -C(=O)- or -S(=O)₂- ;

wherein:

Cy is independently:

C₃₋₂₀carbocyclyl,

C₃₋₂₀heterocyclyl, or

C₅₋₂₀aryl;

and is optionally substituted;

Q¹ is independently:

a covalent bond;

C₁₋₇alkylene; or

C₁₋₇alkylene-X-C₁₋₇alkylene, -X-C₁₋₇alkylene, or C₁₋₇alkylene-X-,

wherein X is -O- or -S-;

and is optionally substituted;

Q² is independently:

C₄₋₈alkylene;

and is optionally substituted;

and has a backbone length of at least 4 atoms;

or:

Q² is independently:

C₅₋₂₀arylene-C₁₋₇alkylene;

~~C₁₋₇alkylene-C₅₋₂₀arylene; or,~~

~~—C₁₋₇alkylene-C₅₋₂₀arylene-C₁₋₇alkylene;~~

and is optionally substituted;

and has a backbone length of at least 4 atoms;

or a pharmaceutically acceptable salt, ~~amide, ester, or ether thereof,~~

provided that Cy is not pyridine, pyrimidine, a bicyclic ring containing one nitrogen atom, or a bicyclic ring containing at least one of a sulfur or oxygen.

Claims 81-173. (Canceled)

174. (New) A compound according to claim 80, wherein the piperazin-1,4-diyl group is unsubstituted or substituted at one or more the 2-, 3-, 5-, and 6-positions with C₁₋₄alkyl.

175. (New) A compound according to claim 174, wherein J¹ is a covalent bond and J² is -C(=O)-.

176. (New) A compound according to claim 174, wherein J^1 is $-C(=O)-$ and J^2 is $-C(=O)-$.

177. (New) A compound according to claim 174, wherein J^1 is a covalent bond and J^2 is $-S(=O)_2-$.

178. (New) A compound according to claim 174, wherein J^1 is $-C(=O)-$ and J^2 is $-S(=O)_2-$.

179. (New) A compound according to claim 174, wherein Q^1 is independently a covalent bond.

180. (New) A compound according to claim 175, wherein Q^1 is independently a covalent bond.

181. (New) A compound according to claim 176, wherein Q^1 is independently a covalent bond.

182. (New) A compound according to claim 177, wherein Q^1 is independently a covalent bond.

183. (New) A compound according to claim 174, wherein Q^1 is independently C_{1-7} alkylene, and is optionally substituted.

184. (New) A compound according to claim 175, wherein Q^1 is independently C_{1-7} alkylene, and is optionally substituted.

185. (New) A compound according to claim 176, wherein Q^1 is independently C_{1-7} alkylene, and is optionally substituted.

186. (New) A compound according to claim 177, wherein Q^1 is independently C_{1-7} alkylene, and is optionally substituted.

187. (New) A compound according to claim 174, wherein Q^1 is independently C_{1-3} alkylene, and is optionally substituted with one or more groups selected from -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -OPr, -Ph, -NH₂, -CONH₂, and =O.

188. (New) A compound according to claim 175, wherein Q^1 is independently C_{1-3} alkylene, and is optionally substituted with one or more groups selected from -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -OPr, -Ph, -NH₂, -CONH₂, and =O.

189. (New) A compound according to claim 176, wherein Q^1 is independently C_{1-3} alkylene, and is optionally substituted with one or more groups selected from -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -OPr, -Ph, -NH₂, -CONH₂, and =O.

190. (New) A compound according to claim 177, wherein Q^1 is independently C_{1-3} alkylene, and is optionally substituted with one or more groups selected from -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -OPr, -Ph, -NH₂, -CONH₂, and =O.

191. (New) A compound according to claim 174, wherein Q^1 is independently C_{1-3} alkylene-X- C_{1-3} alkylene, -X- C_{1-3} alkylene, or C_{1-3} alkylene-X- wherein X is -O- or -S- and is optionally substituted with one or more groups selected from -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -OPr, -Ph, -NH₂, -CONH₂, and =O.

192. (New) A compound according to claim 174, wherein Q^1 is independently C_{1-3} alkylene-X- C_{1-3} alkylene, -X- C_{1-3} alkylene, or C_{1-3} alkylene-X- wherein X is -O- or -S-.

193. (New) A compound according to claim 174, wherein Q^2 is independently C_{4-8} alkylene and is optionally substituted.

194. (New) A compound according to claim 174, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

195. (New) A compound according to claim 179, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

196. (New) A compound according to claim 180, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

197. (New) A compound according to claim 181, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

198. (New) A compound according to claim 187, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

199. (New) A compound according to claim 188, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

200. (New) A compound according to claim 189, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

201. (New) A compound according to claim 192, wherein Q^2 is independently a saturated aliphatic C_{4-8} alkylene group.

202. (New) A compound according to claim 174, wherein Q^2 is independently selected from $-(CH_2)_5-$, $-(CH_2)_6-$, $-(CH_2)_7-$, and $-(CH_2)_8-$.

203. (New) A compound according to claim 174, wherein Q^2 is independently C_{5-20} arylene- C_{1-7} alkylene and is optionally substituted.

204. (New) A compound according to claim 174, wherein Q^2 is independently C_{5-6} arylene- C_{1-7} alkylene and is optionally substituted.

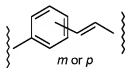
205. (New) A compound according to claim 174, wherein Q^2 , is independently phenylene- C_{1-7} alkylene, C_{1-7} alkylene-phenylene, or C_{1-7} alkylene-phenylene- C_{1-7} alkylene and is optionally substituted.

206. (New) A compound according to claim 174, wherein Q^2 , is independently phenylene-methylene, phenylene-ethylene, or phenylene-ethenylene and is optionally substituted.

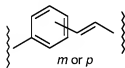
207. (New) A compound according to claim 206, wherein the phenylene linkage is meta.

208. (New) A compound according to claim 206, wherein the phenylene linkage is para.

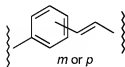
209. (New) A compound according to claim 174, wherein Q^2 , is independently:



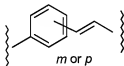
210. (New) A compound according to claim 179, wherein Q^2 , is independently:



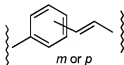
211. (New) A compound according to claim 182, wherein Q^2 , is independently:



212. (New) A compound according to claim 187, wherein Q^2 , is independently:



213. (New) A compound according to claim 190, wherein Q^2 , is independently:



214. (New) A compound according to claim 174, wherein Q^2 has a backbone of at least 5 atoms.

215. (New) A compound according to claim 174, wherein Q^2 has a backbone of at least 6 atoms.

216. (New) A compound according to claim 174, wherein Cy is independently C₅₋₂₀carboaryl or C₅₋₂₀heteroaryl and is optionally substituted.

217. (New) A compound according to claim 174, wherein Cy is independently phenyl, furanyl, pyrrolyl, imidazolyl, pyrazinyl, pyridizinyl, naphthyl, fluorenyl, acridinyl, or carbazolyl; and is optionally substituted.

218. (New) A compound according to claim 174, wherein Cy is independently phenyl or naphthyl; and is optionally substituted.

219. (New) A compound according to claim 174, wherein Cy is independently phenyl and is optionally substituted.

220. (New) A compound according to claim 179, wherein Cy is independently phenyl and is optionally substituted.

221. (New) A compound according to claim 187, wherein Cy is independently phenyl and is optionally substituted.

222. (New) A compound according to claim 194, wherein Cy is independently phenyl and is optionally substituted.

223. (New) A compound according to claim 195, wherein Cy is independently phenyl and is optionally substituted.

224. (New) A compound according to claim 196, wherein Cy is independently phenyl and is optionally substituted.

225. (New) A compound according to claim 197, wherein Cy is independently phenyl and is optionally substituted.

226. (New) A compound according to claim 198, wherein Cy is independently phenyl and is optionally substituted.

227. (New) A compound according to claim 199, wherein Cy is independently phenyl and is optionally substituted.

228. (New) A compound according to claim 200, wherein Cy is independently phenyl and is optionally substituted.

229. (New) A compound according to claim 201, wherein Cy is independently phenyl and is optionally substituted.

230. (New) A compound according to claim 209, wherein Cy is independently phenyl and is optionally substituted.

231. (New) A compound according to claim 210, wherein Cy is independently phenyl and is optionally substituted.

232. (New) A compound according to claim 211, wherein Cy is independently phenyl and is optionally substituted.

233. (New) A compound according to claim 212, wherein Cy is independently phenyl and is optionally substituted.

234. (New) A compound according to claim 213, wherein Cy is independently phenyl and is optionally substituted.

235. (New) A compound according to claim 174, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -C(=O)NH₂, -C(=O)NMe₂, -C(=O)NEt₂, -C(=O)N(iPr)₂, -C(=O)N(CH₂CH₂OH)₂, -C(=O)Me, -C(=O)Et, -C(=O)cHex, -C(=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

236. (New) A compound according to claim 179, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),

-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

237. (New) A compound according to claim 180, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe,

-C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),
-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,

-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

238. (New) A compound according to claim 181, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

239. (New) A compound according to claim 182, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,

-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

240. (New) A compound according to claim 187, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

241. (New) A compound according to claim 188, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),

-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

242. (New) A compound according to claim 189, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe,

-C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),
-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,

-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

243. (New) A compound according to claim 190, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

244. (New) A compound according to claim 195, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,

-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

245. (New) A compound according to claim 196, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

246. (New) A compound according to claim 197, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),

-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

247. (New) A compound according to claim 198, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe,

-C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),
-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,

-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

248. (New) A compound according to claim 199, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

249. (New) A compound according to claim 200, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,

-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

250. (New) A compound according to claim 210, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂, -(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)cHex, -(C=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

251. (New) A compound according to claim 211, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),

-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

252. (New) A compound according to claim 212, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe,

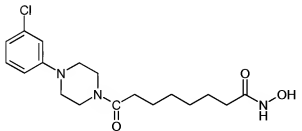
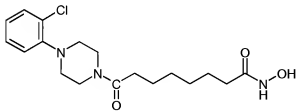
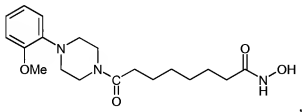
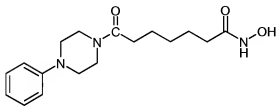
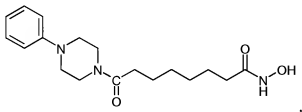
-C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu),
-C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe,
-C(=O)OCH₂CH₂OEt, -(C=O)NH₂, -(C=O)NMe₂, -(C=O)NEt₂, -(C=O)N(iPr)₂,
-(C=O)N(CH₂CH₂OH)₂, -(C=O)Me, -(C=O)Et, -(C=O)-cHex, -(C=O)Ph, -F, -Cl, -Br, -I,
-OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH,
-OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂,
-OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et,
-nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe,
-CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me,

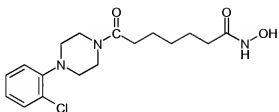
-Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂,
-SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

253. (New) A compound according to claim 213, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)OEt, -C(=O)O(Pr), -C(=O)O(iPr), -C(=O)O(nBu), -C(=O)O(sBu), -C(=O)O(iBu), -C(=O)O(tBu), -C(=O)O(nPe), -C(=O)OCH₂CH₂OH, -C(=O)OCH₂CH₂OMe, -C(=O)OCH₂CH₂OEt, -C(=O)NH₂, -C(=O)NMe₂, -C(=O)NEt₂, -C(=O)N(iPr)₂, -C(=O)N(CH₂CH₂OH)₂, -C(=O)Me, -C(=O)Et, -C(=O)-cHex, -C(=O)Ph, -F, -Cl, -Br, -I, -OH, -OMe, -OEt, -O(iPr), -O(tBu), -OPh, -OCF₃, -OCH₂CF₃, -OCH₂CH₂OH, -OCH₂CH₂OMe, -OCH₂CH₂OEt, -OCH₂CH₂NH₂, -OCH₂CH₂NMe₂, -OCH₂CH₂N(iPr)₂, -OPh, -OPh-Me, -OPh-OH, -OPh-OMe, -OPh-F, -OPh-Cl, -OPh-Br, -OPh-I, -Me, -Et, -nPr, -iPr, -nBu, -iBu, -sBu, -tBu, -nPe, -CF₃, -CH₂CF₃, -CH₂CH₂OH, -CH₂CH₂OMe, -CH₂CH₂OEt, -CH₂CH₂NH₂, -CH₂CH₂NMe₂, -CH₂CH₂N(iPr)₂, -CH₂-Ph, -Ph, -Ph-Me, -Ph-OH, -Ph-OMe, -Ph-F, -Ph-Cl, -Ph-Br, -Ph-I, -SO₂Me, -SO₂Et, -SO₂Ph, -SO₂NH₂, -SO₂NMe₂, -SO₂NEt₂, -NMe₂, -NEt₂, morpholino, -NO₂, and -CN.

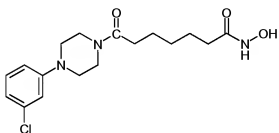
254. (New) A compound according to claim 174, wherein Cy is independently phenyl and is optionally substituted with one or more groups selected from -C(=O)OMe, -C(=O)O(Pr), -C(=O)NHMe, -C(=O)Et, -C(=O)Ph, -OCH₂CH₂OH, -OMe, -OPh, -nPr, -iPr, -CF₃, -CH₂CH₂OH, -CH₂CH₂NMe₂, -Ph, -Ph-F, -Ph-Cl, -SO₂Me, -SO₂Me₂, -NMe₂, -F, -Cl, -Me, -Et, -OMe, -OEt, -CH₂-Ph, and -O-CH₂-Ph.

255. (New) A compound according to claim 80, selected from the following compounds, and pharmaceutically acceptable salts thereof:

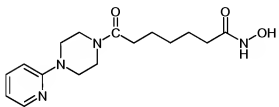




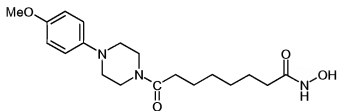
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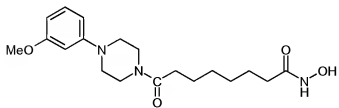
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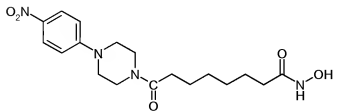
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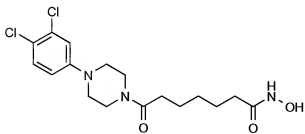
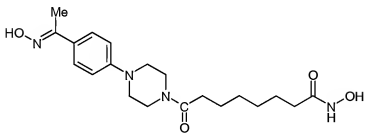
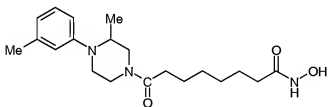
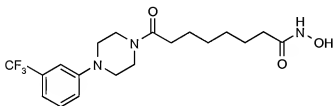
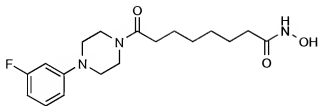
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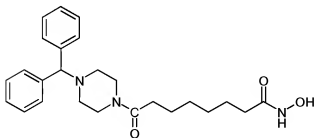
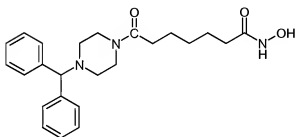
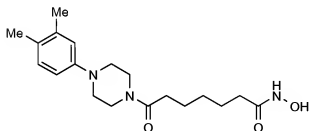
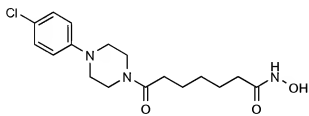
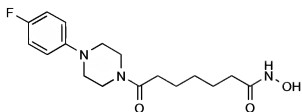


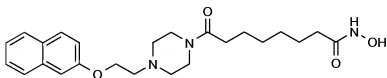
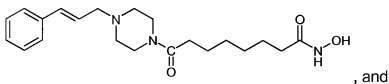
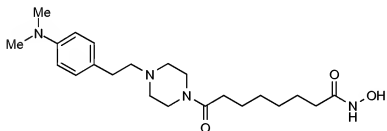
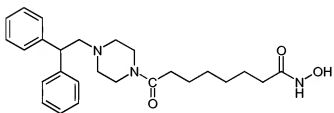
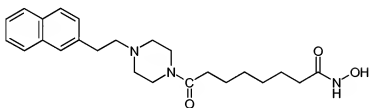
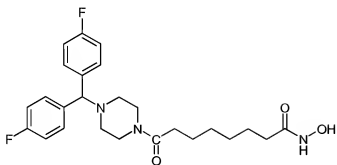
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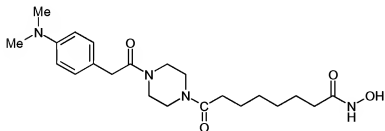
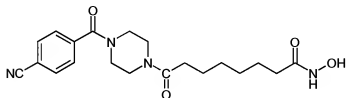
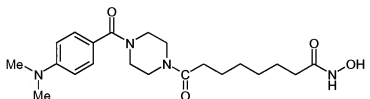
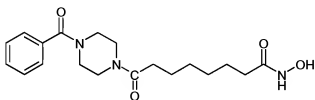
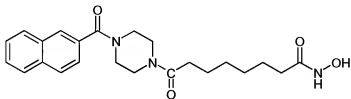
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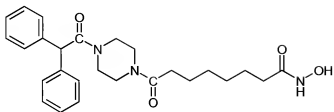




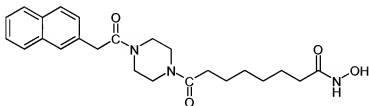


256. (New) A compound according to claim 80, selected from the following compounds, and pharmaceutically acceptable salts thereof:

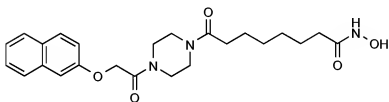




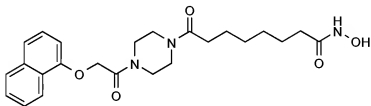
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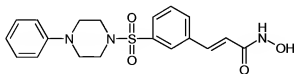


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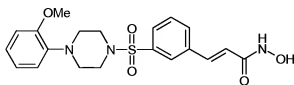


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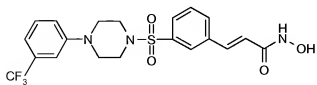
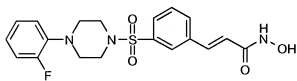
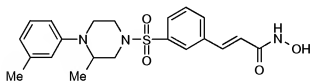
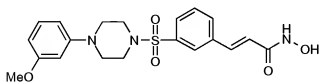
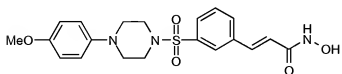
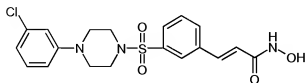
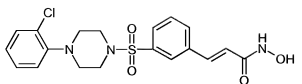
257. (New) A compound according to claim 80, selected from the following compounds, and pharmaceutically acceptable salts thereof:

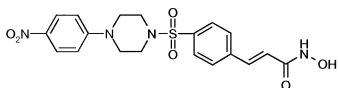
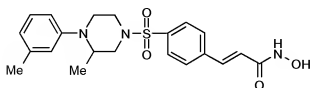
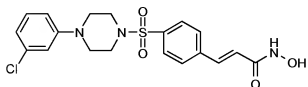
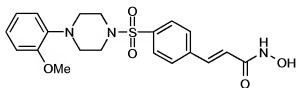
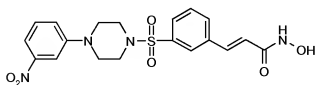
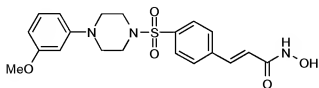
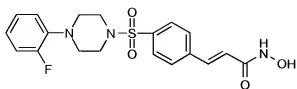


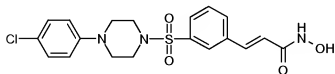
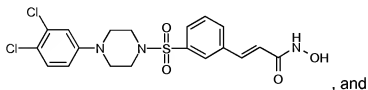
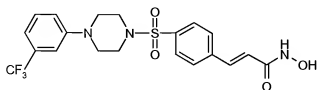
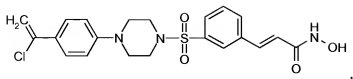
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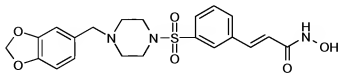
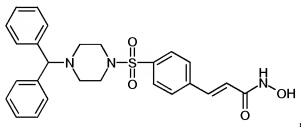
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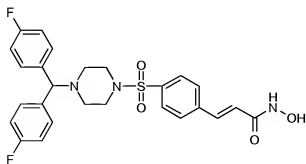
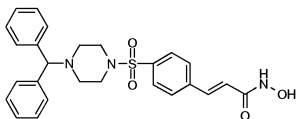
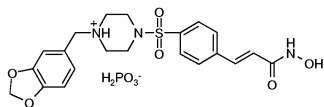
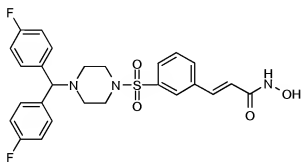
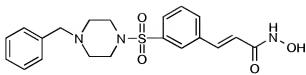


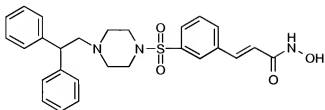




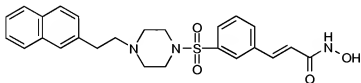
258. (New) A compound according to claim 80, selected from the following compounds, and pharmaceutically acceptable salts thereof:



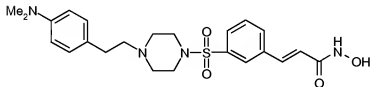




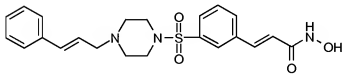
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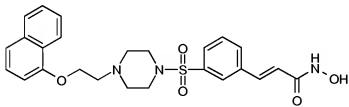
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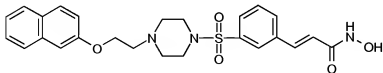
,



,



, and



.

259. (New) A composition comprising a compound according to claim 80 and a pharmaceutically acceptable carrier.